

wherein:

R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub> and R<sub>6</sub>, which are the same or different, are chosen from the group consisting of: H, C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, canphane, adamantane, phenyl, biphenyl, naphthyl, or naphthyl-C<sub>1-8</sub>;

R<sub>5</sub> is chosen from the group consisting of: H, C<sub>1-8</sub> alkyl, C<sub>1-8</sub>alkyl-phenyl, biphenyl, naphthyl, COOR, CN or phenyl,;

X is chosen from the group consisting of: O, C(=O)R, COOR, NO<sub>2</sub>, and CONNR', wherein R and R' are as above defined;

Q is chosen from the group consisting of single-bond, C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, canphane, adamantane, CO, CONR, and NR, where R is as above defined;

W is chosen from the group consisting of H, C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, canphane, adamantane, trifluoromethyl, C<sub>1-8</sub> alkoxy, C<sub>1-8</sub> alkoxy-C<sub>1-8</sub> alkyl, phenyl, biphenyl, naphthyl-C<sub>1-8</sub> alkyl, phenyl, biphenyl, naphthyl, phenyloxy, biphenyloxy, naphthyloxy, phenylamino, biphenylamino, naphthylamino, C<sub>1-8</sub> alkyl-carbonyl, phenylcarbonyl, biphenylcarbonyl, naphthylcarbonyl, phenylcarboxyl, biphenylcarboxyl, naphthylcarboxyl, phenylcarboxamide, biphenylcarboxamide, naphthylcarboxamide, halogen, CN, NRR', C<sub>1-8</sub> alkylamino;

n is an integer comprised between 1 and 4;

the symbol ----- means that the corresponding bonds a, b, c, d, e, f, g, h and i are single or double bonds, with the proviso that when b or f are a double bond, the group R<sub>5</sub> is absent; their pharmaceutically acceptable salts and esters.

2. (twice amended) A benzo(c)quinolizine compound of formula (1) according to Claim 1, wherein R<sub>5</sub> = H, C<sub>1-8</sub> alkyl-phenyl, biphenyl, naphthyl;

X = O, COOH;

Q = single bond, CO, CONR, NR, wherein R

is chosen from the group consisting of H, C<sub>1-8</sub> alkyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, canphane, adamantane, phenyl, biphenyl, naphthyl or naphthyl-C<sub>1-8</sub>alkyl;

W = H, F, Cl, Br, Me, t-butyl, C<sub>1-8</sub>alkoxy, 2,5-dimethylhexyl, trifluoromethyl, 2,5-(di-trifluoromethyl)-phenyl, 4-methyloxy-phenyl, phenyl, phenyl-C<sub>1-8</sub>alkyl, C<sub>1-8</sub>alkylcarbonyl, phenylcarbonyl;

n = 1 and 2;

R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub> and R<sub>5</sub> = H, Me, CN, phenyl, COOR, CONRR', C(=O)R, wherein R and R' are the same or different and are chosen from the group consisting of H, C<sub>1-8</sub> alkyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, canphane, adamantane, phenyl, biphenyl, naphthyl or naphthyl-C<sub>1-8</sub>.

3. (amended) A benzo[c]quinolizine compounds according to Claim 1 of the formula:

2,3,4,4a,5,6,6a,7,8,9,10,10a-dodecahydro-(1H)-benzo[c]quinolizin-3-one;

8-chloro-2,3,4,4a,5,6,6a,7,8,9,10,10a-dodecahydro-(1H)-benzo[c]quinolizin-3-one;

2,3,4,4a,5,6,6a,7,8,9,10,10a-dodecahydro-8-methyl-(1H)-benzo[c]quinolizin-3-one;

2,3,4,4a,5,6,6a,7,8,9,10,10a-dodecahydro-4-methyl-(1H)-benzo[c]quinolizin-3-one;

2,3,4,4a,5,6,6a,7,8,9,10,10a-dodecahydro-1-methyl-(1H)-benzo[c]quinolizin-3-one;

2,3,4,4a,5,6,6a,7,8,9,10,10a-decahydro-(1H)-benzo[c]quinolizin-3-one;

8-chloro-2,3,4,4a,5,6,6a,7,8,9,10,10a-decahydro-(1H)-benzo[c]quinolizin-3-one;

2,3,4,4a,5,6,6a,7,8,9,10,10a-decahydro-8-methyl-(1H)-benzo[c]quinolizin-3-one;

2,3,5,6,6a,7,8,9,10,10a-decahydro-4-methyl-(1H)-

benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-1-methyl-(1H)-benzo[c]quinoliz-  
zin-3-one;  
(4a $\alpha$ ,6a $\beta$ ,10a $\alpha$ )-3,4,5,6,6a,7,8,9,10,10a-decahydro-(4aH)-  
benzo[c]quinoliz-in-3-one;  
[(4a $\alpha$ ,6a $\beta$ ,10a $\alpha$ )-3,4,5,6,6a,7,8,9,10,10a-decahydro-(4aH)-  
benzo[c]quinoliz-in-3-one;]  
3,4,5,6,6a,7,8,9,10,10a-decahydro-(1H)-benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-(4aH)-  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-8-methyl-(4aH)-  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-4-methyl-(4aH)-  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-1-methyl-(4aH)-  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-4-methyl-(1H)-  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-4,8-dimethyl-(1H)-  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-1-methyl-(1H)-  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-1,4-dimethyl-(1H)-  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-4-methyl-(4aH)-  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-4,8-dimethyl-(4aH)-  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-1-methyl-(4aH)-  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-1,8-dimethyl-(4aH)-  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-5-methyl-(1H)-  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-5-methyl-(1H)-  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-5,8-dimethyl-(1H)-

benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-4,5-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-1,5-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-5-methyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-5-methyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-5,8-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-4,5-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-1,5-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-4,5-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-4,5,8-trimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-1,5-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-1,4,5-trimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-4,5-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-4,5,8-trimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-1,5-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-1,5,8-trimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-6-methyl-(1H) -  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-6-methyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-4,6-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;

E<sup>1</sup>  
6<sup>4</sup>  
cont

2,3,5,6,6a,7,8,9,10,10a-decahydro-1,6-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-6,8-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-6,8-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-4,6-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-1,6-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-4,6-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-4,6,8-trimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-1,6-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-1,4,6-trimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-4,6-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-4,6,8-trimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-1,6-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-1,6,8-trimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-5,6-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-5,6-dimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-4,5,6-trimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-1,5,6-trimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-5,6-dimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-5,6-dimethyl-(4aH) -

benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-5,6,8-trimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-4,5,6-trimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-1,5,6-trimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-1,5,6-trimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-4,5,6-trimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-4,5,6,8-tetramethyl-(1H) -  
benzo[c]quinolizin-3-one;  
8-chloro-2,3,5,6,6a,7,8,9,10,10a-decahydro-1,5,6-trimethyl-(1H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,6a,7,8,9,10,10a-decahydro-1,4,5,6-tetramethyl-(1H) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-4,5,6-trimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-4,5,6,8-tetramethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-1,5,6-trimethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
3,4,5,6,6a,7,8,9,10,10a-decahydro-1,5,6,8-tetramethyl-(4aH) -  
benzo[c]quinolizin-3-one;  
5,6,6a,7,8,9,10,10a-octahydro-(3H) -benzo[c]quinolizin-3-one;  
8-chloro-5,6,6a,7,8,9,10,10a-octahydro-(3H) -benzo[c]quinolizin-3-  
one;  
5,6,6a,7,8,9,10,10a-octahydro-8-methyl-(3H) -benzo[c]quinolizin-3-  
one;  
5,6,6a,7,8,9,10,10a-octahydro-4-methyl-(3H) -benzo[c]quinolizin-3-  
one;  
8-chloro-5,6,6a,7,8,9,10,10a-octahydro-4-methyl-(3H) -  
benzo[c]quinolizin-3-one;  
5,6,6a,7,8,9,10,10a-octahydro-4,8-dimethyl-(3H) -  
benzo[c]quinolizin-3-one;  
2,3,5,6,7,8,9,10-octahydro-(1H) -benzo[c]quinolizin-3-one;

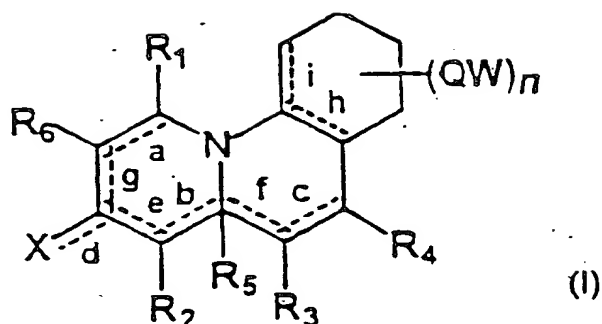
8-chloro-2,3,5,6,7,8,9,10-octahydro-(1H)-benzo[c]quinolizin-3-one;  
 2,3,5,6,7,8,9,10-octahydro-8-methyl-(1H)-benzo[c]quinolizin-3-one;  
 2,3,5,6,6a,7,8,9-octahydro-(1H)-benzo[c]quinolizin-3-one;  
 8-chloro-2,3,5,6,6a,7,8,9-octahydro-(1H)-benzo[c]quinolizin-3-one;  
 2,3,5,6,6a,7,8,9-octahydro-8-methyl-(1H)-benzo[c]quinolizin-3-one;  
 4a-benzyl-3,4,5,6,6a,7,8,9,10,10a-decahydro-(4aH)-benzo[c]quinolizin-3-one;  
 4a-benzyl-8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-(4aH)-benzo[c]quinolizin-3-one;  
 4a-benzyl-3,4,5,6,6a,7,8,9,10,10a-decahydro-8-methyl-(4aH)-benzo[c]quinolizin-3-one;  
 4a-benzyl-3,4,5,6,6a,7,8,9,10,10a-decahydro-4-methyl-(4aH)-benzo[c]quinolizin-3-one;  
 4a-benzyl-3,4,5,6,6a,7,8,9,10,10a-decahydro-1-methyl-(4aH)-benzo[c]quinolizin-3-one;  
 3,4,5,6,6a,7,8,9,10,10a-decahydro-4a-(4-pyridyl)methyl-(4aH)-benzo[c]quinolizin-3-one;  
 8-chloro-3,4,5,6,6a,7,8,9,10,10a-decahydro-4a-(4-pyridyl)-methyl-(4aH)-benzo[c]quinolizin-3-one;  
 3,4,5,6,6a,7,8,9,10,10a-decahydro-8-methyl-4a-(4-pyridyl)methyl-(4aH)-benzo[c]quinolizin-3-one;  
 3,4,5,6,6a,7,8,9,10,10a-decahydro-4-methyl-4a-(4-pyridyl)methyl-(4aH)-benzo[c]quinolizin-3-one[;].  
 3,4,5,6,6a,7,8,9,10,10a-decahydro-1-methyl-4a-(4-pyridyl)methyl-(4aH)-benzo[c]quinolizin-3-one;

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27. (amended) A method for the inhibition of 5 $\alpha$  reductase-I and/or 5 $\alpha$  reductase-II iso-enzymes as defined in claim 13 where the pathology is selected from the group consisting of acne, baldness, prostatic cancer and prostatic hypertrophy in men and hirsutism in women.

28. (amended) A fully and partially reduced benzo(c)quinolizine

compound of formula (1):



wherein:

*E2*  
*Cont.*  
R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub> and R<sub>6</sub>, which are the same or different, are chosen from the group consisting of: H, C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, camphane, adamantane, phenyl, biphenyl or naphthyl;

R<sub>5</sub> is chosen from the group consisting of: H, C<sub>1-8</sub> alkyl, C<sub>1-8</sub>alkyl-phenyl, biphenyl, naphthyl, COOR, CN or phenyl;

X is chosen from the group consisting of: O, C(=O)R, COOR, NO<sub>2</sub>, and CONNR', wherein R and R' are as above defined;

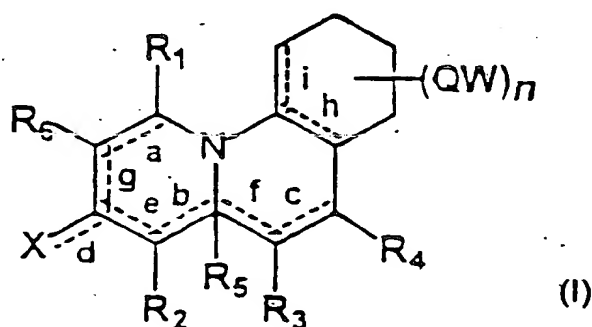
Q is chosen from the group consisting of single-bond, C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, camphane, adamantane, CO, CONR, and NR, where R is as above defined;

W is chosen from the group consisting of H, C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, camphane, adamantane, trifluoromethyl, C<sub>1-8</sub> alkoxy, C<sub>1-8</sub> alkoxy-C<sub>1-8</sub> alkyl, phenyl, biphenyl, naphthyl-C<sub>1-8</sub> alkyl, phenyl, biphenyl, naphthyl, phenyloxy, biphenyloxy, naphthyloxy, phenylamino, biphenylamino, naphthylamino, C<sub>1-8</sub> alkyl-carbonyl, phenylcarbonyl, biphenylcarbonyl, naphthylcarbonyl, phenylcarboxyl, biphenylcarboxyl, naphthylcarboxyl, phenylcarboxamide, biphenylcarboxamide, naphthylcarboxamide, halogen, CN, NRR' where R and R' are as above defined;



n is an integer comprised between 1 and 4;  
the symbol ----- means that the corresponding bonds a, b, c, d, e, f, g, h and i are single or double bonds, with the proviso that when b or f are a double bond, the group R<sub>5</sub> is absent; their pharmaceutically acceptable salts and esters.

29. (new) A fully and partially reduced benzo(c)quinolizine compound of formula (1):



wherein:

R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub> and R<sub>5</sub>, which are the same or different, are chosen from the group consisting of: H, C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, camphane, adamantane, phenyl, biphenyl or naphthyl;

R<sub>5</sub> is chosen from the group consisting of: H, C<sub>1-8</sub> alkyl, C<sub>1-8</sub>alkyl-phenyl, biphenyl, naphthyl, COOR, CN, phenyl;

X is chosen from the group consisting of: O, C(=O)R, COOR, NO<sub>2</sub>, and CONNR', wherein R and R' are as above defined;

Q is chosen from the group consisting of single-bond, C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, camphane, adamantane, CO, CONR, and NR, where R is as above defined;

W is chosen from the group consisting of H, C<sub>1-8</sub> alkyl, C<sub>2-8</sub> alkenyl, C<sub>2-8</sub> alkynyl, cyclopropane, cyclobutane, cyclopentane, cyclohexane, cycloheptane, cyclooctane, norbornane, camphane, adamantane, trifluoromethyl, C<sub>1-8</sub> alkoxy, C<sub>1-8</sub> alkoxy-C<sub>1-8</sub> alkyl,